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L5 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2007 ACS ON STN ACCESSION NUMBER: 2002.565410 CAPLUS DOCUMENT NUMBER: 137:201298

TITLE:

Preparation of substituted isoxazolines as anti-depressants
Andres-Gil, Jose Ignacio, Pernandez-Gadea, Francisco Javier, Alcazar-Vaca, Manuel Jesus; Cid-Nunez, Jose Mariar, Pastor-Fernandez, Joaquin; Megens, Antonius Adrianus Hendrikus Petrus; Heylen, Godelieve Irma Christine Maria; Langlois, Xavier Jean Michel; Bakker, Margareha Henrica Maria; Steckler, Thomas Horst INVENTOR (S):

Janssen Pharmaceutica N.V., Belg. PCT Int. Appl., 100 pp. CODEN: PIXXD2 PATENT ASSIGNEE(S): SOURCE: DOCUMENT TYPE:

Patent English 1 FAMILY ACC. NUM. COUNT: PATENT INFORMATION: LANGUAGE:

APPLICATION NO. KIND PATENT NO.

20020213 <--CA, CH, CN, GD, GE, GH, I.C, I.K, I.R, NZ, OM, PH, A, TR, TT, TZ, M. AT' AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
M. AT' AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
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SS, FI, FR, GB, GR, IE, IT, LU, MC, NL, FT, SE, TR,
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AI 20020829 CA 2002-2437505 20020213 <-AI 20030934 AU 2002-2437505 20020213 <-AI 2003094 AU 2002-243717
BI 20031210 EP 2002-712909 20020213
BI 20060821 CK, IT, LI, LU, NL, SE, MC, PT, 466333335 W: 2002066484
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OTHER SOURCE(S): GI	MARPA.	MARPAT 137:201298					

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Title compds. I [wherein X = CH2, NR7, S or 0; R7 = H, (un)substituted alkyl, Ph. Ph alkyl, etc.; R1 and R2 independently = H, OH, CN, halo, 0802H, (un)substituted Ph, phenylalkyl, alkoxy, etc.; or R1 and R2 may be taken together to form a bivalent radical selected from -CH2CH2O., -OCH2CH2., -OCH2CH2 and -OCH2CH2O.; m = 1-4; Y = (un)substituted piperidyl or piperaxyl radical and R3 represents an (un)substituted aromatic homocyclic or heterocyclic ring system including a partially or completely hydrogenated hydrocarbon chain of maximum 6 atoms long with which the ring æ

Erich Leese <12/04/2007>

10/513699

system is attached to the Y radical and which may contain one or more heteroatoms selected from the group of 0. N and S], a process for their preparation, pharmaceutical compins. comprising them and their use as a medicine for treating anxiety disorders and disorders of body weight are disclosed. Thus, II was prepared in 60% yield by reaction of III with N-(3-phenyl-2-propenyl)-piperazine. III was prepared by substitution of Me 4-bromo-2-butenoate with 2-hydroxy-4,5-dimethoxybenzaldehyde with subsequent condensation with methanesulfont of the invention have surprisingly been shown to have a scrotonin (5-HT) reuptake inhibitor activity in combination with addni. a2-adrenoceptor antagonist activity and show a strong anti-depressant activity without often also at the haza and hack size (but the 1-HT ransporter site of more than 50% (piC50) at a test concentration ranging between 10-6 M and 10-9 M in a concentration dependent manner. The invention also relates to novel combination of substituted isoxazolines derive having anti-depressant activity and/or anxiolytic activity and/or body weight control activity with antidepressants, anxiolytic activity and/or antipasychotics to improve efficacy and/or anxiolytic activity and/or 42321-33-69-48 452321-31-69-48 452321-39-69-68 452321-31-99-89 452321-31-99-99 452322-19-9

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RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
(Reactant or reagent) (intermediate; preparation and pharmaceutical activity of substituted isoxazolines as anti-depressants)

452321-67-4 CAPLUS Z Z

3H-[1]Benzopyrano[4,3-c]isoxazole-3-carboxylic acid, 3a,4-dihydro-7,8-dimethoxy-, methyl ester, (3R,3aS)-rel- (9Cl) (CA INDEX NAME)

Relative stereochemistry.

3H-[1] Benzopyrano [4,3-c] isoxazole-3-methanol, 3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) 452321-69-6 CAPLUS Z Z

Relative stereochemistry

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S S

452321-71-0 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole-3-methanol, 3a,4-dihydro-7,8-dimethoxy-, methanesulfonate (ester), (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452321-73-2 CAPLUS CN 3H-[1]Benzopyrano[4,3-c]isoxazole-3-methanol, 3a,4-dihydro-7-hydroxy-8-methoxy-, α-methanesulfonate, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

S &

452311-75-4 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-(1-piperazinylmethyl)-, (3K,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007>

Erich Leese

10/513699

452321-77-6 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole-3-methanol, 3a,4-dihydro-7,8-dimethoxy-, methanesulfonate (ester), (3R,3aS)-rel-(+)- (9CI) (CA INDEX NAME) S &

Rotation (+). Absolute stereochemistry unknown.

RN 452321-80-1 CAPLUS CN 3H-[1]Benzopyrano[4,3-c]isoxazole-3-methanol, 3a,4-dihydro-7,8-dimethoxy-, methanesulfonate (ester), (3R,3aS)-rel-(-)- (9CI) (CA INDEX NAME)

Rotation (-). Absolute stereochemistry unknown.

452321-62-3 CAPLUS
1-Piperazincearboxylic acid, 4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-, 1,1-dimethylethyl ester, rel[9CI] (CA INDEX NAME) Z Z

Relative stereochemistry.

452321-85-6 CAPLUS
2-Propanone, 1-[4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]-, rel- (9CI)
(CA INDEX NAME) Z Z

Relative stereochemistry.

452321-87-8 CAPLUS Piperazine, 1-[[(3x,3aS)-3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-4-[(3,4-dihydro-2-naphthalenyl)carbonyl]-, rel-(9Cl) (CA INDEX NAME) **Z** Z

Relative stereochemistry.

452321-89-0 CAPLUS
hosphonium, [2-(1(3R,)as])-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-(1)soxazol-3-yl]methyl]-1-piperazinyl]ethyl]triphenyl-,
bromide, rel- (9CI) (CA INDEX NAME) Z Z

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10/513699

Relative stereochemistry.

Br-

452321-91-4 CAPLUS
Piperazine, 1-[([3R,3aS]-3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-6]isoxazol-3-yl]methyl]-4-(3-iodo-2-methyl-1-oxo-2-propenyl)-, rel- [9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry unknown.

452121-93-6 CAPLUS
452121-93-6 (A-(3-iodo-2-methyl-2-properyl)-1-piperazinyl]methyl]-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) **3** 3

Relative stereochemistry. Double bond geometry unknown.

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RN 452321-95-8 CAPLUS
CN 1-Piperazineethanol, 4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[(4,3-c)isoxazol-3-yl]methyl]- a-methyl-, rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

NN 452321-97-0 CAPLUS

CN 1,2-Propanediol, 3-[4-[([3R,3aS)-3a,4-dlhydro-7,8-dlmethoxy-3H[1]benzopyrano[(4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]-, rel- (9CI)

(CA INDEX NAME)

Relative stereochemistry.

RN 452321-99-2 CAPLUS
CN 1-Piperazineacetaldehyde, 4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

10/513699

RN 452322-05-3 CAPLUS
CN Isoxazolo[4,3-c]quinoline-3-carboxylic acid, 3,3a,4,5-tetrahydro-5(trifluoroacetyl)-, methyl ester, (3R,3aR)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452322-07-5 CAPLUS CN Isoxazolo[4,3-c]quinoline-3-methanol, 3,3a,4,5-tetrahydro-, (3R,3aR)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

Relative stereochemistry.

RN 452322-19-9 CAPLUS
CN 1-Piperazinecarboxylic acid, 4-[[(3R,3aR)-3,3a,4,5-tetrahydro-7,8-dimethoxy-5-(trifluoroacetyl)isoxazolo[4,3-c]quinolin-3-yl]methyl]-,
1,1-dimethylethyl ester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

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452322-21-3 CAPLUS

1-Piperazinecarboxylic acid, 4-[[(3R,3aR)-3,3a,4,5-tetrahydro-7,8-dimethoxyisoxazolo[4,3-c]quinolin-3-yl]methyl]-, 1,1-dimethylethyl ester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

452122-23-5 CAPLUS Isoxazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-(1-piperazinylmethyl)-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry.

Z Z

452132-29-1 CAPLUS
1-Piperidinecarboxylic acid, 4-[[[[3R,3aS]-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]amino]methyl]-,
1,1-dimethylethyl ester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007>

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10/513699

Z 23

452322-30-4 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole-3-methanamine, 3a,4-dihydro-7,8-dimethoxy-N-(4-piperidinylmethyl)-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

H

452322-32-6D, resin bound 452323-46-5D, resin bound
RL: FCT (Reactant); RACT (Reactant or reagent)
('preparation and pharmaceutical activity of substituted isoxazolines as anti-depressants)
452322-32-6 CAPUS
31-11]Benzopyrano(4,3-c)isoxazole-3-methanol, 3a,4-dihydro-7,8-dimethoxy-, methanesulfonate (ester) (9CI) (CA INDEX NAME) Z Z

Z Z

452323-46-5 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[2-(diphenylphosphino]ethyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA

<12/04/2007>

INDEX NAME)

Relative stereochemistry.

H

452313-12-5P 452313-68-7P 452313-71-2P
452313-80-3P 452313-87-5P 452316-78-8P
RL: PAC (Pharmacological activity); PUR (Purification or recovery); SPN
(Synthetic preparation); THU (Therapeutic use); BIOL (Biological study);
PREP (Preparation); USES (Uses)
(target compound, preparation and pharmaceutical activity of substituted isoxacolines as anti-depressants)
452313-32-5 CAPLUS
3H-[1]Benzopyrano[4]-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-pipperazinyl]methyl]-, dihydrochloride,
(3R,3aS)-rel-(+)- (9CI)

S S

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

●2 HCl

Z Z

452313-68-7 CAPLUS Isoxazole(4,3-c]quinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-2-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel-(+)-(9CI) (CA INDEX NAME)

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

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10/513699

452313-71-2 CAPLUS
Isoxazolo(4,2-clquinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[(4-[(2E)-2-19opthyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel-(-)-(9CI) (CA INDEX NAME) **3** 3

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

452313-80-3 CAPLUS
31[1]Beracpyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, dihydrochloride, [3R,3aS]-rel-(+)- (9CI) (CA INDEX NAME) S S

Rotation (+). Absolute stereochemistry unknown.

●2 HCl

452313-82-5 CAPLUS

88

<12/04/2007>

3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, dihydrochloride,
(3R,3aS)-rel-(-)- (9CI) (CA INDEX NAME) S

Rotation (-). Absolute stereochemistry unknown.

●2 HCl

452316-78-8 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-[1]Benzopyrano[1]-1-piperaziny]methyl]-, dihydrochloride, (3R,3aS)-rel-(-)- (9CI) (CA INDEX NAME) Z Z

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

D2 HCl

H

RL: FAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic RL: FAC (Pharmacological activity); BIOL (Biological study); PREP (Preparation); TRU (Therapeutic use); BIOL (Biological study); PREP (Erage compound; preparation and pharmaceutical activity of substituted isoxazolines as anti-depressants)
452313-59-6 CAPLUS
ISOxazolo(4,3-c)quinoline, 3,3a,4,5-tetrahydro-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) Z Z

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<12/04/2007>

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Relative stereochemistry.

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452314-83-9P 452314-86-2P 452314-89-5P 452314-98-6P 452318-92-0P 452314-89-5P 452314-89-5P 452314-98-6P 452315-10-4P 452315-04-7P 452315-10-1P 452315-13-1P 452315-19-4P 452315-19-1P 452315-19-4P 452315-19-1P 452315-19-4P 452315-19-1P 452315-19-4P 452315-19-1P 452315-19-4P 452315-19-1P 452315-1P-4P 452315-1P-4P 452315-4P-1P 45231
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452313-98-3P 452313-91-6P 452313-93-8P
452314-08-8P 452314-10-1P 452314-0-5P
452314-16-8P 452314-11-3P 452314-14-6P
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452314-31-7P 452314-34-0P 452314-37-3P
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452316-24-4P 452316-27-P 452316-30-2P
452316-33-5P 452316-36-8P 452316-39-1P
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452317-20-3P 452317-22-5P 452317-24-7P
452317-26-9P 452317-28-1P 452317-30-5P
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452316-51-7P 452316-53-9P
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452117-32-7P 452117-34-9P 452317-36-1P
452317-38-3P 452317-40-7P 452317-42-9P
452317-50-9P 452317-6-19 452317-6-9P
452317-50-9P 452317-5-7P 452317-6-1P
452317-7-10-PP 452317-67-PP 452317-6-1P
452317-7-11-4P 452317-7-6-PP 452317-7-9-PP
452317-7-11-4P 452317-8-7P 452317-8-9P
452317-9-2P 452317-8-7P 452317-9-9P
452317-9-2P 452318-1P-8-7P 452317-9-9P
452317-9-4P 452318-1P-8-7P 452318-7-9P
452318-15-9P 452318-1P-9P
452318-15-9P 452318-1P-9P
452318-15-9P 452318-1P-7P
452318-15-9P 452318-1P-7P
452318-1P-4P 452318-1P-PP
452318-1P-4P-PP
452318-1P-4P 452318-1P-PP
452318-1P-4P-PP
452319-1P-4P-PP
45
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(target compound, preparation and pharmaceutical activity of substituted isoxazolines as anti-depressants) Z Z

452313-36-9 CAPLOS
31-11]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CAINDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452313-40-5 CAPLUS 2

Erich Leese <12/04/2007>

10/513699

3H-[1]Benzopyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) Z

Relative stereochemistry. Double bond geometry as shown.

452313-43-8 CAPLUS
31-[1] Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(6-quinolinylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry.

452313-46-1 CAPLUS
3H-[1] Benzogytzano[4,3-c] isoxazole, 3-[[4-[(2E)-3-(4-fluorophenyl)-23H-[1] Benzogenyl]-1-piperazinyl methyl]-3a, 4-dihydro-7,8-dimethoxy-,
[3R,3aS]-rel-(-)- [9CI) (CA INDEX NAME) 2 2

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

Erich Leese <12/04/2007>

452313-50-7 CAPLUS ₹ 3

3H-[1]Benzopyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, acetate (ester),
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

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452313-54-1 CAPLUS

4.1]-Renzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenzy-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

Z Z

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

<12/04/2007>

10/513699

452313-61-0 CAPLUS Isoxazolo[4,3-c]quinoline, 5-acetyl-3,3a,4,5-tetrahydro-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry.

452313-65-4 CAPLUS Isoxazalol(4,2elquinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) 2 2

Relative stereochemistry. Double bond geometry as shown.

452313-74-5 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(3-chlorophenyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME) **3 3**

<12/04/2007>

RN 452313-77-8 CAPLUS
CN 34-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

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Relative stereochemistry.

RN 452313-85-8 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methy]-3-(3-thieny])-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

owie boild geometry as shown.

RN 452313-88-1 CAPLUS
CN 3H-[1]Benzopyrano(4,3-c)isoxazole, 3-[[4-[(4-chlorophenyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

<12/04/2007> Erich Leese

10/513699

Relative stereochemistry.

RN 452313-91-6 CAPLUS
CN 34-(1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(2-methylphenyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

RN 452313-93-8 CAPLUS
CN 3+[1]Benzopyranol4,3-c]isoxazole, 3-[[4-[(3,4-dihydro-2naphthaleny])methyl]-1-piperaxinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452313-98-3 CAPLUS CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-3-(2-pyridinyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-

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(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Z Z

452314-01-1 CAPLUS
3H-(11Benzogyano(4,3-c)isoxazole, 3-[(4-[(2E)-3-(4-fluorophenyl)-2-methyl2-propenyl)-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

452314-05-5 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(3-phenyl-2-butenyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX

Relative stereochemistry. Double bond geometry unknown.

Erich Leese <12/04/2007>

10/513699

452314-08-8 CAPLUS

1-Piperazineethanamine, 4-[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-α-methyl-N-phenyl-, rel-(9CI) (CA INDEX NAME) C. Z.

Relative stereochemistry.

452314-11:3 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[2-(3-fluorophenoxy)propyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME) Z Z

452314-14-6 CAPLUS
3H-[1]Benzopyrano[4, 3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(2-naphthanylanethyl)-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel-(9CI) (CA INDEX NAME) S S

Relative stereochemistry.

Erich Leese <12/04/2007>

•2 HCl

Z Z

452314-16-8 CAPLUS
3H-11] Henzopyrano[4,3-c] isoxazole-3-methanamine, 3a,4-dihydro-7,8-dimethox-N-[1-(2-naphthalenylmethyl)-4-piperidinyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452314-18-0 CAPLUS
CN 3H-[1] Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(phenylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

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10/513699

RN 452314-20-4 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[[2-(rifluoromethyl]phenyl]methyl]-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME)

RN 452314-23-7 CAPLUS
CN 31-[1] Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[[3-(trifluoromethy1)pheny1]methy1]-1-piperaziny1]methy1] (9CI) (CA INDEX NAME)

RN 452314-26-0 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[[4-[Krifluoromethyl]phenyl]methyl]-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME)

RN 452314-29-3 CAPLUS CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[[2-

<12/04/2007>

(1.1,2,2.tetrafluoroethyl)phenyl]methyl]-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME)

RN 452314-31-7 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(2-bromophenyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME)

RN 452314-34-0 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[{4-[(3-bromophenyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME)

RN 452314-37-3 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(4-bromophenyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME)

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Relative stereochemistry.

RN 452314-43-1 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(4-f]luorophenyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, dihydrochloride,
(3K,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

•2 HCl

RN 452314-46-4 CAPLUS

<12/04/2007>

ű CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(2-fluorophenyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI)
INDEX NAME)

Relative stereochemistry.

452314-49-7 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(4-nitrophenyl)methyl]-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME) Z Z

452314-52-2 CAPLUS
Benzenamine, 4-[[4-[(3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]methyl]-N.N-dimethyl- (GA INDEX NAME) Z Z

Z Z

452314-55-5 CAPLUS
Acetamide, N-[4-[[4-[(3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl)methyl]-1-piperazinyl]methyl]phenyl]- (9CI) (CA INDEX

<12/04/2007>

Erich Leese

NAME) 10/513699

452314-57-7 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(2-ethoxyphenyl)methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME) Z Z

452314-60-2 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(3-phenoxyphenyl)methyl]-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME) S K

452314-62-4 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(4phenoxyphenyl)methyl]-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME) Z Z

<12/04/2007>

452314-65-7 CAPLUS

1-Propanamine, 3-[4-[(4-[(3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl)methyl]-1-piperazinyl]methyl]phenoxy]-N,N-dimethyl- (9CI)
(CA INDEX NAME) £ 5

452114-68-0 CAPLUS Benzoic acid, 4-[[4-[(3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl)methyl]-l-piperazinyl]methyl]-, methyl ester (9CI) (CA INDEX NAME) Z Z

452314-71-5 CAPLUS
Benzonitrile, 3-[[4-[(3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl)methyl]-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME) Z Z

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<12/04/2007>

452314-74-8 CAPLUS
Benzonitrile, 4-[[4-[(3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl)methyl]-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME) **3** 3

452314-77-1 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-([1,1'-biphenyl]-3-ylmethyl)-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CAINDEX NAME) Z Z

Relative stereochemistry.

452314-80-6 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(2,4-dimethoxyphenyl)methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME) ₩ 53

RN 452314-83-9 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(3,4-dimethoxyphenyl)methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- [9CI) (CA INDEX NAME)

RN 452314-86-2 CAPLUS
CN 3H-[i]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(2,4-dichlorophenyl]methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry.

RN 452314-89-5 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(3,4-dichlorophenyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry.

<12/04/2007>

Erich Leese

10/513699

RN 452314-92-0 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(4-bromo-2-fluorophenyl)methyl]1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

RN 452314-95-3 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(1-methyl-1H-pyrrol-2-yl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

RN 452314-98-6 CAPLUS
CN 3H-11]Benzépyranolt,3-c}isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2-phenyl-1H-imidazol-4-yl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

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Relative stereochemistry.

RN 452315-01-4 CAPLUS CN 3H-[1] Benzopyrano[4,3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[2-thiazolylmethyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-04-7 CAPLUS
CN 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2-grangut-4-thiazolyl)methyl]-1-piperazinyl]methyl]-1-piperazinyl]methyl]-(2-grangut-1

Relative stereochemistry.

452115-07-0 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(2-furanylmethyl)-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA

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<12/04/2007> Erich Leese

10/513699

INDEX NAME)

Relative stereochemistry.

RN 452315-10-5 CAPLUS
CN 34-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(5-bromo-2-furanyl)methyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CL) (CA
INDEX NAME)

Relative stereochemistry.

RN 452315-13-8 CAPLUS
CN 34-[1] Benzopyrano[4,3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(5-methyl-2-furanyl)methyl]-- (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-16-1 CAPLUS CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(3-furanylmethyl)-1-

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piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-19-4 CAPLUS
CN 3H-[1] Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(3-phenyl-1,2,4-oxadiazol-5-yl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-22-9 CAPLUS
CN 3H-[1]senzopyrand(4,3-c)isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(5-phenyl-1,3,4-oxadiazol-2-yl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-24-1 CAPLUS

<12/04/2007>

Erich Leese

10/513699

CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(2-thienylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-27-4 CAPLUS
CN 34-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(5-bromo-2-thienyl)methyl]-1-INDEX NAME)
INDEX NAME)

Relative stereochemistry.

RN 452315-30-9 CAPLUS
CN 34-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(4-bromo-2-thieny1)methy1]-1piperazinyl]methy1]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

RN 452315-33-2 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(3-thenylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX

Relative stereochemistry.

RN 452315-36-5 CAPLUS
CN 3H-[1]Benzopyrano(4,3-c]isoxazole, 3-[[(3R,5S)-3,5-dimethyl-4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-38-7 CAPLUS
CN 3H-[1] Benzopyrano[4,3-c] isoxazole, 3-[[(2R,5S)-2,5-dimethyl-4-(2-naphthalenylmethyl)-1-piperazinyl] methyl] -3a,4-dihydro-7,8-dimethoxy-, (3S,3aR)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007> Brich Leese

10/513699

RN 452315-40-1 CAPLUS
CN 3H-[1] Benzopyrano[4,3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[3-methyl-4-(2-naphthalenylmethyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-42-3 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(1-naphthalenylmethyl)-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME)

RN 452315-44-5 CAPLUS
CN 3H-(11Benzopayranol4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(1-naphthalenylmethyl)-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007>

•2 HCl

452315-46-7 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(1-methoxy-2-naphthalenyl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry.

S S

452315-48-9 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[[6methoxy-2-naphthalenyl]methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

RN 452315-51-4 CAPLUS

<12/04/2007>

Erich Leese

10/513699

3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(3-methoxy-2-naphthalenyl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

452315-52-5 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(7-methyl-2-naphthalenyl)methyl]-1-piperazinyl]methyl}-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) S S

Relative stereochemistry.

452315-55-8 CAPLUS
#1[]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[[6-methy1-2-naphthaleny1]methy1]-1-piperaziny1]methy1]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry.

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RN 452315-58-1 CAPLUS
CN 3H-[1]8encopyrano(4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(3-methyl-2-naphthalenyl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

RN 452315-61-6 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(7-fluoro-2-naphthalenyl)methyl]1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

RN 452315-63-8 CAPLUS
CN 34-[1] Renzopyrano[4,3-c] isoxazole, 3-[[4-[(6-fluoro-2-naphthalenyl)methyl]1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-; (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

10/513699

RN 452315-66-1 CAPLUS
CN 3+-[1]Renzoprano[4,3-c]isoxazole, 3-[[4-[(5-fluoro-2-naphthalenyl)methyl]1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

RN 452315-70-7 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(5-chloro-2-naphthalenyl)methyl]i-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

RN 452315-73-0 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(1-bromo-2-naphthalenyl)methyl]1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

<12/04/2007>

RN 452315-76-3 CAPLUS
CN 3H-[1]Benzapyranol(4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(5-nitro-2-naphthalenyl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME)

Relative stereochemistry.

RN 452315-79-6 CAPLUS
CN 1-Naphthalenamine, 6-[[4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]methyl]-, rel(9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-82-1 CAPLUS
CN 1-Nephthalenamine, 6-[[4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]methyl]-N.Ndimethyl-, rel- (9CI) (CA INDEX NAME)

<12/04/2007>

Erich Leese

10/513699

Relative stereochemistry.

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RN 452315-85-4 CAPLUS
CN 3H-[1]Berrapyranof4,3-c]isoxazole, 3-[[4-[[6,7-difluoro-2naphthalenyl]methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-87-6 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(5,7-difluoro-2-naphthalenyl)methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-90-1 CAPLUS

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CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(7,8-difluoro-2-naphthalenyl)methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-92-3 CAPLUS
CN 3H-[1]Berrapyranof(4,3-c]isoxazole, 3-[[4-[[5-bromo-8-methoxy-2-naphthalenty]]methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452315-94-5 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-3-[[4-(1H-indol-2-ylmethyl)-1-piperazinyl]methyl]-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CAINDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

10/513699

RN 452315-97-8 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-3-[[4-(1H-indol-3ylmethyl)-1-piperazinyl]methyl]-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry.

RN 452316-00-6 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(1H-benzimidazol-2-ylmethyl)-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, dihydrochloride,
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

2 HCl

RN 452316-03-9 CAPLUS
CN 34-[1]Benzoprano[4,3-c]isoxazole, 3-[[4-(2-benzofuranylmethyl)-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, dihydrochloride, (3R,3aS)-rel- (91) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

2 HC1

RN 452316-06-2 CAPLUS
CN #1-[1] Benzopyranol(4,3-c] isoxazole, 3-[[4-(benzo[b]thien-2-ylmethyl)-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, dihydrochloride,
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

•2 HCl

RN 452316-09-5 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(benzo[b]thien-3-ylmethyl)-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

10/513699

RN 452316-12-0 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-{[4-(2-quinolinylmethyl)-1-piperazinyl]methyl]- (9CI) (CA INDEX NAME)

RN 452316-15-3 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(3-quinolinylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452316-18-6 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(4-quinolinylmethyl)-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007>

●2 HCl

RN 452316-21-1 CAPLUS CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(2-phenylethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452316-24-4 CAPLUS
CN 3H-[1]Benzogyzano[4,3-c]isoxazole, 3-[[4-[2-(1,3-benzodioxol-5-yl)ethyl]-1-piperazinyl]methyl]-3a.4-dihydror-7,8-dimethoxy-, dihydrochloride, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

•2 HCl

<12/04/2007> Erich Leese

10/513699

RN 452316-27-7 CAPLUS
CN 34-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[2-(2-nghthaleny])ethyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

●2 HCl

RN 452316-30-2 CAPLUS
CN 34-11Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[2-(1-naphthaleny])ethyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

2 HCl

RN 452316-33-5 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(3-phenylpropyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

RN 452316-36-8 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-3-[[4-(1H-inden-2ylmethyl)-1-piperazinyl]methyl]-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry.

RN 452316-39-1 CAPLUS
CN 34-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(1H-benzimidazol-2-ylmethyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452316-42-6 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[2-(4-fluorophenoxy)ethyl]-1-piperaziny]]methyl]-3-4-dihydro-7,8-dimethoxy-, dihydrochloride, (3R,3aS)-rel- (9C) RNDEX NAME)

Relative stereochemistry.

<12/04/2007>

Erich Leese

10/513699

•2 HCl

RN 452316-45-9 CAPLUS
CN 3# [1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[2-(1,3-benzodioxol-5yloxy]ethyl]--:piperazinyl]methyl].3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452316-48-2 CAPLUS
CN 1H-Inden-1-one, 5-[2-[4-[[3R,3aS]-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyranol(4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]ethoxy]-2,3-dihydro-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452316-51-7 CAPLUS

<12/04/2007>

CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[2-[(2,3-dihydro-5-benzofuranyl)oxy]ethyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

452316-53-9 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[2-(3-f]luorophenoxy)propyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CAINDEN NAME)

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Relative stereochemistry.

RN 452316-55-1 CAPLUS
CN 1(2H)-Benzogotrannon, 6-[2-[4-[[(3R,3a5)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzogoyrano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]-1-methylethoxy], rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

10/513699

RN 452316-58-4 CAPLUS
CN 3H-[1] Bergopyrano[4,3-c] isoxazole, 3-[[4-[[(2R)-6-fluoro-3,4,4a,8a-tetrahydro-2H-1-benzopyran-2-yJ]methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452316-64-2 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(1,3-benzodioxol-2-ylmethyl)-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry.

RN 452316-66-4 CAPLUS
CN 1-Propanone, 3-(4-1(18r,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]-1-phenyl-, rel(9CI) (CA INDEX NAME)

Relative stereochemistry.

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1-Piperazinepropanol, 4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]- a-phenyl-, rel- (9CI) (CA INDEX NAME) 452316-69-7 CAPLUS Z Z

Relative stereochemistry.

3 3

452316-72-2 CAPLUS
1-Propanone, 3-14-[[(3R, 3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1] Denzopyrano[4,3-c] isoxazol-3-yl] methyl]-1-piperazinyl]-1-phenyl-,
oxime, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry unknown.

3 3

452316-75-5 CAPLUS
4F1[]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CAINDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

452316-81-3 CAPLUS Ph-11]Benzopyrano(4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(3-ph-nyl-2-propenyl)-1-piperazinyl]methyl]-, (3K,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452316-84-6 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(22)-3-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452316-87-9 CAPLUS
3H-[1] Benzoyyrano[4,3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3]-(1] Benzoyphenyl] -2-propenyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z 23

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Relative stereochemistry. Double bond geometry as shown.

●2 HCl

452316-89-1 CAPLUS
4F1]lBencopyrano[4, 3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(3-methoxyphenyl)-2-propenyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel- (9C) (Ca INDEX NAME) Z 3

Relative stereochemistry. Double bond geometry as shown.

•2 HCl

Z Z

452316-91-5 CAPLUS

**IllBencapyranot (4.3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(4-methoxyphenyl)-2-propenyl)-1-piperazinyl]methyl]-, dihydrochloride, (18,3aS)-rel- (901) (A INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Erich Leese

<12/04/2007>

10/513699

•2 HCl

452316-93-7 CAPLUS
3H-[1] Benzopyrano[4,3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(4-phenoxyphenyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry unknown.

452316-95-9 CAPLUS
3H-[1]Benzoppyrano[4, 3-c] isoxazole, 3-[[4-[(2E)-3-(2-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

RN 452316-97-1 CAPLUS
CN 3H-[1]Benzopyrano[4, 3-c]isoxazole, 3-[[4-[(2E)-3-(3-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452316-99-3 CAPLUS
3H-[11Benzopyrano|4,3-c]isoxazole, 3-[[4-[3-(3-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CL) (CAINDEX NAME)

Z Z

Relative stereochemistry. Double bond geometry unknown.

452317-02-1 CAPLUS
3H-[11.Benzopyranol4,3-clisoxazole, 3-[[4-[(2E)-3-(4-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,dihydrochloride, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

●2 HCl

452317-04-3 CAPLUS
Aphth[1,2-c]isoxazole, 3-[[4-[(2E)-3-(4-fluorophenyl)-2-propenyl]-1-piperazinyl]methyll-3,3a,4,5-tetrahydro-7,8-dimethoxy-, (3R,3aS)-rel-(+)-(9Cl) (CA INDEX NAME) Z Z

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

452317-06-5 CAPLUS
3H-[1] Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(2-chlorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA Z Z

Relative stereochemistry. Double bond geometry unknown.

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3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(3-chlorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) 452317-08-7 CAPLUS Z Z

Relative stereochemistry. Double bond geometry unknown.

452317-10-1 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(4-chlorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3as)-rel- (9Cl) (CAINDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

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452317-12-3 CAPLUS
4-(1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(2-bromophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

Erich Leese <12/04/2007>

10/513699

452317-14-5 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(3-bromopheny])-2-propeny]]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA **3** 33

Relative stereochemistry. Double bond geometry unknown.

452317-16-7 CAPLUS

4-1]Benzopyrano(4,3-clisoxazole, 3-[4-[3-(4-bromophenyl)-2-propenyl]-1-piperazinolyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CAINDEX NAME) 2 E

Relative stereochemistry. Double bond geometry unknown.

452317-18-9 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-[2-(trifluoromethyl)phenyl]-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) 2 Z

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Relative stereochemistry. Double bond geometry unknown.

Z Z

452317-20-3 CAPLUS

1-11]Bennopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-[3-[3-[1:1]luoromethy]]pheny]]-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry unknown.

Z Z

Relative stereochemistry. Double bond geometry unknown.

Erich Leese <12/04/2007>

10/513699

452317-24-7 CAPLUS
3H-[1] Benzogyyano[4, 3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-[2-(1,1,2,2-terrafluoroethyl) phenyl]-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry unknown.

452317-26-9 CAPLUS
Benzoic acid, 4-13-(4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1)benzopyzano[4,3-c]:soxazol-3-yl]methyl]-1-piperazinyl]-1-propenyl]-,
methyl ester, rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452317-28-1 CAPLUS

Benzoic acid, 3-[3-[4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c];soxazol-3-yl]methyl]-1-piperazinyl]-1-propenyl]-,
methyl ester, rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

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452317-30-5 CAPLUS
4F2317-30-5 CAPLUS
4F1[]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(3-nitrophenal)]-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry unknown.

452317-32-7 CAPLUS

4F1118enzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(2-nitropheny])-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) £ 5

Relative stereochemistry. Double bond geometry unknown.

452317-34-9 CAPLUS
3H-[1] Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(3-methylphenyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

<12/04/2007>

Erich Leese

10/513699

Relative stereochemistry. Double bond geometry unknown.

452317-36-1 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(3-phenoxyphenyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) C KN

Relative stereochemistry. Double bond geometry unknown.

452317-38-3 CAPLUS
Benzonitrile, 2-[3-[4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c]soxazol-3-yl]methyl]-1-piperazinyl]-1-propenyl]-,
rel- (9CI) (CA INDE NAME) **3** 3

Relative stereochemistry. Double bond geometry unknown.

Erich Leese <12/04/2007>

Z Z

452317-40-7 CAPLUS
Benzonitzile, 3-13-[4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyzano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]-1-propenyl]-,
rel- (9Cl) (CA INDEX RAME)

Relative stereochemistry. Double bond geometry unknown.

3 3

452317-42-9 CAPLUS
Benzonitrile, 4-[3-44-[(13K,3aS)-3a,4-dihydro-7,8-dimethoxy-3H[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]-1-propenyl]-,
rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.. Double bond geometry unknown.

452317-44-1 CAPLUS
452317-44-1 CAPLUS
4.1-1.1-1.1-2.2.2.2.2.2.3.2.3.3.4-dihydro-7.8-dimethoxy-, (3R,3aS)-rel-9c() (A. INDEX NAME) Z 5

Relative stereochemistry. Double bond geometry unknown.

Erich Leese

<12/04/2007>

10/513699

452317-46-3 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(3,4-difluoropheny])-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CAINDEX NAME) S S

<u>5</u> 452317-48-5 CAPLUS
3H-[I]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-[2,4-difluorophenyl]-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI)
INDEX NAME) S &

452317-50-9 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(2,3-difluorophenyl]-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CAINDEX NAME) Z Z

Erich Leese

452317-52-1 CAPLUS
4F1[lBencopyrano[4,3-c]isoxazole, 3-[[4-[3-(2,6-difluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-[9CI] (CA INDEX NAME) £ 5

Relative stereochemistry. Double bond geometry unknown.

452117-54-3 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(2,5-difluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) **3 3**

Relative stereochemistry. Double bond geometry unknown.

452317-56-5 CAPLUS Z.

<12/04/2007>

10/513699

CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(3,5-difluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry unknown.

3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(2,6-dichloropheny1)-2-propeny1]-1-piperaziny1]methy1]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) 452317-58-7 CAPLUS ₹ 8

Relative stereochemistry. Double bond geometry unknown.

452117-60-1 CAPLUS

4-11]Benzopyrano(4,3-c]isoxazole, 3-[(4-[3-(2-chloro-4-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9Cl) (CA INDEX NAME) S 5

<12/04/2007>

S S

Relative stereochemistry. Double bond geometry unknown.

452317-67-8 CAPLUS

LilBenzopyrano[4,3-c] isoxazole, 3-[4-[(2E)-3-(2,3-dihydro-1H-inden-5-yl)-2-propenyl]-1-piperazinyl]nethyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452317-69-0 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(2,3-dihydro-1,4-benzodioxin-6-2 Z

Erich Leese <12/04/2007>

10/513699

yl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry unknown.

452317-71-4 CAPLUS
3H-[1]Benzoyyzano[4,3-c]isoxazole, 3-[[4-[3-(9H-fluoren-2-yl)-2-propenyl]1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) Z 23

Relative stereochemistry. Double bond geometry unknown.

452317-73-6 CAPLUS
3H-[11Benzopranol4,3-c]isoxazole, 3-[[4-[3-(2-dibenzofurany1)-2-propenyl]1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry unknown.

452317-76-9 CAPLUS
4-11]senzopyranol(43-c1)soxazole, 3-[(4-[(2E)-3-(1,3-benzodioxol-5-y1)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452317-79-2 CAPLUS
3H-[1]Bencopyranol(4,3-c]isoxazole, 3-[[4-[(2E)-3-(2,3-dihydro-5-benzofuranyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-(3R,3aS)-rel-(9CI) [(9CI INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452317-82-7 CAPLUS
4-11/Bencopyranol(4)-clisoxazole, 3-[[4-[3-(4,5-dimethoxy-2-nitrophenyl)-2-propenyl]-1-piperazinyl]methyl]-a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (AR INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

<12/04/2007>

10/513699

452317-84-9 CAPLUS
3H-[11Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(3-thienyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452317-86-1 CAPLUS
4F11]sencopyrano(4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(3-thienyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(+)- (9CI) (CA INDEX NAME) Z Z

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

452317-89-4 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(3-thienyl)-2-propenyl]-1-piperaçinyl]methyl]-, (3K,3aS)-rel-(-)- (9CI) Z Z

Erich Leese

(CA INDEX NAME)

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

Z Z

452317-92-9 CAPLUS
4-11]emcopyranol(4,3-c]isoxazole, 3-[[4-[3-(4-bromo-2-thienyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-9Cl) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry unknown.

Z Z

Br

452317-94-1 CAPLUS
3H-[1] Benzogyano[4, 3-6] isoxazole, 3-[[4-[3-(5-bromo-2-thienyl)-23H-[1] Benzogyano[4, 3-6] isoxazole, 3-[[4-[3-(5-bromo-2-thienyl)-2propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-,
[9CI] (CA INDEX NAME)

Relative stereochemistry. Double bond geometry unknown.

Erich Leese <12/04/2007>

10/513699

452317-96-3 CAPLUS

4-118enzopyrano(4,3-c]isoxazole, 3-[[4-[3-(2-furanyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452317-99-6 CAPLUS
3H-[1] Benzopyrano[4,3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(3-pyridinyl) -2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452318-02-4 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(2-pyridinyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452318-04-6 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-[5-(4-chlorophenyl]-2-furanyl]-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, [3R,3aS]-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452318-07-9 CAPLUS
3H-[11Benzopyrano|4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(2-naphthalenyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452318-09-1 CAPLUS
4F11]Rencapyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(1-naphthalenyl)-2-propenyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel- (9CI) (CA INDEX NAME) ₹ 3

<12/04/2007>

Erich Leese

10/513699

Relative stereochemistry. Double bond geometry as shown.

•2 HCl

452318-11-5 CAPLUS
3H-(11Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(4-quinoliny)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452318-13-7 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[3-(2-quinoliny])-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CAIDEX NAME) S S

Relative stereochemistry. Double bond geometry unknown.

Z Z

452318-15-9 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[3-(9-anthracenyl)-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME)

PAGE 1-A

PAGE 2-A

Z Z

452318-18-2 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethòxy-3-[[4-[(2E)-1-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)

Erich Leese <12/04/2007>

10/513699

(CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452318-20-6 CAPLUS
3H-[11Benzopyrano[4, 3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(+)-(9CI) (CA INDEX NAME) Z Z

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

452318-22-8 CAPLUS
3H-[1] Benzopyzano[4, 3-c] isoxazole, la,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, dihydrochloride, [3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

●2 HCl

3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(-)-(9CI) (CA INDEX NAME) 452318-24-0 CAPLUS S S

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

452318-27-3 CAPLUS
4-4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-8-4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-8-4-1]-1]-4-4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-, (2E)-2-butenedioate (1:2) (9CI) (CA INDEX NAME) S S

δ

CRN 452318-26-2 CMF C27 H33 N3 O4

Relative stereochemistry. Double bond geometry as shown.

Erich Leese

<12/04/2007>

10/513699

N £ CRN 110-17-8 CMF C4 H4 O4

Double bond geometry as shown.

452318-30-8 CAPLUS
3H:(11 Benropyranol4,)-c]isoxazole, 3-[[(2R,5S)-2,5-dimethyl-4-[(2E)-2-methyl-3-phenyl-2-properyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) 2 S

Relative stereochemistry. Double bond geometry as shown.

£ 5

452318-32-0 CAPLUS
3H-[1] Benzopyrano(4,3-c]isoxazole, 3-{[(3R,5S)-3,5-dimethyl-4-{(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9C1) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452318-34-2 CAPLUS
3H-[1]Benzopyranole, 3-c]isoxazole, 3-([(2R,6S)-2,6-dimethyl-4-[(2E)-2-methyl-3-phenyl-2-propenyl-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452318-16-4 CAPLUS

H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2Z)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S 53

Relative stereochemistry. Double bond geometry as shown.

452318-38-6 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(2-methyl-3-phenyl-2-propenyl]-1-plperazinyl]methyl]- (9CI) (CA INDEX NAME) $\overset{C}{\mathbb{R}}\overset{K}{\mathbb{R}}$

Erich Leese

<12/04/2007>

10/513699

452318-41-1 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(2,3-diphenyl-2-propenyl)-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy- (9CI) (CA INDEX NAME) S S

$$\begin{array}{c|c} & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ & &$$

452318-43-3 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(2-[1,1'-biphenyl]-4-yl-3-phenyl-2-propenyl)-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry unknown.

Erich Leese

Z 5

452318-45-5 CAPLUS

**IlBenzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[2-(2-naphthalenyl)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9C1) (CA_INDEX_NAME)

Relative stereochemistry. Double bond geometry unknown.

452318-47-7 CAPLUS

**IlBenzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[2-[phenylmethylene]butyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry unknown.

452318-49-9 CAPLUS
4F1]Bencopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-(phenylmethylne)heptyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z 3

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

452318-52-4 CAPLUS
3H-[1]Benzopyrano[4, 3-C] isoxazole, 3-[[4-[(22]-2-bxomo-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) £ 5

Relative stereochemistry. Double bond geometry as shown.

452318-54-6 CAPLUS
3H-[11Benzogyrano[4, 3-C]isoxazole, 3-[[4-[(22)-2-chloro-3-phenyl-2-propenyl-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-goll) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452318-57-9 CAPLUS
3H-[1]Benzopyrano[4, 3-c]isoxazole, 3-[[4-[(22)-2-fluoro-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9C1) (CA INDEX NAME) Z Z

<12/04/2007>

Relative stereochemistry. Double bond geometry as shown.

452318-60-4 CAPLUS
1-Piperazinepropanenitrile, 4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H-1)-benzopyzane(4,3-c)isoxazol-3-yl]methyl] - a-(phenylmethylene)-, (a2)-rel- (9Cl) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

Z Z

452318-63-7 CAPLUS
4F1[]Bencopyrano[4,3-c]isoxazole, 3-[[4-[3-[4-(1,1-dimethylethyl)]phenyl]-2-methyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry unknown.

Erich Leese <12/04/2007>

10/513699

452318-65-9 CAPLUS
3H-[1] Benzopyrano[4, 3-c] isoxazole, 3-[[4-(3-[1,1'-biphenyl]-4-yl-2-methyl-2-propnyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CL) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry unknown.

452318-67-1 CAPLUS
3H-[1]Benzopyrano[4,13-C]isoxazole, 3-[[4-[[22]-3-(4-fluorophenyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S 53

Relative stereochemistry. Double bond geometry as shown.

452318-69-3 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(2Z)-3-(3-fluorophenyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

452318-71-7 CAPLUS
3H-(1]Benzogyano(4,3-c)isoxazole, 3-[[4-[(2E)-3-(3-fluorophenyl)-2-methyl-3-flBenzogyarano(4,2-fl)methyl-3-4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9Cl) (CA INDEX NAME) £ 5

Relative stereochemistry. Double bond geometry as shown.

452318-73-9 CAPLUS
3H-[1]Benzogyrano[4,3-c]isoxazole, 3-[[4-[(2E)-3-(4-fluorophenyl)-2-methyl2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452318-75-1 CAPLUS
3H-[1]Benzopyrano[4, 3-c]isoxazole, 3-[[4-[(2E)-3-(4-fluorophenyl)-2-methyl2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel-(-)-(9CI) (CA INDEX NAME) Z Z

Erich Leese <12/04/2007>

10/513699

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

2 Z

452318-77-3 CAPLUS
3H-[11Benzogyrano[4, 3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3]-(2-methoxyphenyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (GA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452318-79-5 CAPLUS
3H-111Benzogyrano[4, 3-c]isoxazole, 3-[[4-[(2E)-3-(2-chloro-6-fluorophenyl)-2-enethyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

452318-81-9 CAPLUS Z Z

3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(2Z)-3-(2,5-difluorophenyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452318-83-1 CAPLUS
41.[1]eenzopyrano[4,3-c]isoxazole, 3-[4-[(2E)-3-(2,5-difluorophenyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

3H-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-[(2Z)-3-(2,3-difluorophenyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl}-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) 452318-85-3 CAPLUS 23

Relative stereochemistry. Double bond geometry as shown.

Erich Leese

<12/04/2007>

10/513699

452318-87-5 CAPLUS

##[1]Benzogyrano[4,3-c]isoxazole, 3-[[4-[(2E)-3-(2,3-difluorophenyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452318-89-7 CAPLUS

[1] Benzopyrano[4,3-c] isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[2-maphthaleny])-2-propeny]]-1-piperazinyl]methyl]-, [3R,3aS)-rel- (9CI) (CA INDEX NAME) S. 25

Relative stereochemistry. Double bond geometry unknown.

452318-91-1 CAPLUS
31-[1]Benzopyrano[4,3-c]isoxazole, 3-[[4-(2H-1-benzopyran-3-ylmethyl)-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CAINDEX RAME) Z Z

Erich Leese

Relative stereochemistry.

452318-93-3 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenyl-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452318-95-5 CAPLUS
4F1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenyl-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(+)- (9CI) (CA INDEX NAME) S S

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

RN 452318-97-7 CAPLUS

<12/04/2007>

Erich Leese

10/513699

CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)3-phenyl-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(-)- (9CI) (CA
INDEX NAME)

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

S S

452318-99-9 CAPLUS Phenol, 3-[(18]-3-[4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-1-piperazinyl]-1-methyl-1-propenyll-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452319-01-6 CAPLUS
3H-[11Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3+1]1Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-10-2-buteny]]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

S S

452319-03-8 CAPLUS
4-11]Bencopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-3-(2-thienyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Z Z

452319-05-0 CAPLUS
4F1]Bencapyrano(43-c]isoxazole, 3-[(4-[(2E)-3-(2,5-dimethyl-3-thienyl)-2-methyl-2-propenyl]-1-piperazinyllmethyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Z Z

452319-07-2 CAPLUS #/Ll/Benzopyranol(4)-clisoxazole, 3-[(4-[(2E)-3-(2-furanyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9Cl) (CA INDEX NAME)

<12/04/2007>

Erich Leese

10/513699

Relative stereochemistry. Double bond geometry as shown.

S &

452319-09-4 CAPLUS
3H-[1]Benzopyrano[4, 3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methy-3-(3-pyridinyl)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452319-11-8 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-2-methyl-3-(2-methyl-4-thiazolyJ)-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z 5

Relative stereochemistry. Double bond geometry as shown.

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RN 452319-13-0 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2-phenyl-1-cyclopenten-1-yl)methyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452319-15-2 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-(3-phenyl-2-propynyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452319-17-4 CAPLUS
CN 3H-[1]Benzopyrano(4,3-c]isoxazole, 3-[[4-[3-[2-(4-fluorophenyl)-1,3-dioxolan-2-yl]propyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

10/513699

RN 452319-20-9 CAPLUS
CN 3+[1]Renzopyrano[4,3-c]isoxazole, 3-[[4-[3-(4-fluorophenoxy)propyl]-1piperaziny]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry.

RN 452319-22-1 CAPLUS
CN 3H-(1]Benzopyrano(4,3-c)isoxazole, 3-[[4-[(2E)-4-(2-fluorophenyl)-2-methyl-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,
(3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

RN 452319-24-3 CAPLUS CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E,4E)-5-phenyl-2,4-pentadienyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-

Erich Leese

(CA INDEX NAME) (BCI) Relative stereochemistry. Double bond geometry as shown.

Z Z

452319-25-4 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452319-27-6 CAPLUS
enzoda caid, 4-[[(18.3-1-4-[[(3R,3aS)-3a,4-dihydro-3H-[1]benzopyrano[4,3-c]isoxazol-3-yllmethyl]-1-piperazinyl]-1-propenyl]-, methyl ester, rel-(9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Z Z

452319-29-8 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 7-fluoro-3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3K,3aS)-rel- (9CI)

<12/04/2007>

10/513699

(CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452319-31-2 CADLUS 3H-[1]Benzopyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) £ 5

Relative stereochemistry. Double bond geometry as shown.

452319-33-4 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S &

Relative stereochemistry. Double bond geometry as shown.

452319-35-6 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7-methoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) 2 Z

Relative stereochemistry.

<12/04/2007>

Double bond geometry as shown.

452319-37-8 CAPLUS

**IlBenzopyrano(4,3-c]isoxazole, 3a,4-dihydro-7-methoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rei-(+, (9Cl) (CA INDEX NAME) Z Z

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

•2 HCl

452319-39-0 CAPLUS
31-(1)Benzopyranol(4,3-c]isoxazole, 3a,4-dihydro-7-methoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel-(-)- (9Cl) (CA INDEX NAME) S &

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

●2 HC1

<12/04/2007>

10/513699

452319-41-4 CAPLUS

**I.!]Bencopyrano[4,3-c]isoxazole, 3a,4-dihydro-7-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

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45230-86-4P 45230-8-6P 452320-90-0P 452320-90-92-2P 452220-94-4P 452320-96-6P 45231-0-96-6P 45231-0-96-6P 45231-10-9P 45231-10-7P 45231-10-1P 452321-10-7P 452321-10-1P 452321-10-P 452321-10-P 452321-10-P 452321-11-9P 452321-21-0P 452321-31-2P 452321-33-4P 452321-33-4P 452321-35-6P 452321-33-4P
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       452319-47-0P
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452319-43-6P 452319-45-8P
452319-49-2P 452319-51-6P
452319-55-0P 452319-57-2P
452319-61-8P 452319-63-0P
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452320-68-2P 452320-70-6P
452320-74-0P 452320-76-2P
452320-80-8P 452320-82-0P
H
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RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(target compound, preparation and pharmaceutical activity of substituted
isoxazolines as anti-depressants)
452119-43-6 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7-(2-methoxyethoxy)-3-[[4-([2E)-2-methy1-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel(9CI) (CA INDEX NAME)

5 5

Relative stereochemistry. Double bond geometry as shown.

452319-45-8 CAPLUS
3H-[1] Benzopyranol (4,3-c) isoxazole, 3a,4-dihydro-7-[(2-methoxyelnexy) -3-[[4-([2E]-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

3 S

452319-47-0 CAPLUS
3H-[1]Benzogyzano[4,3-c]isoxazole, 3a,4-dihydro-7-[(2-methoxyethoxy)methoxy]-1-[4-[(2E)-2-methy]-3-phenyl-3-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

<12/04/2007>

Erich Leese

10/513699

452319-49-2 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a.4-dihydro-7-[(2-methoxyethoxy)methoxy]-3-[[4-[(2E)-2-methy]-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

●2 HCl

452319-51-6 CAPLUS
Carbanic acid, ethyl., (3R,3aS)-3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]:3H-[l]benzopyrano[4,3-G]isoxazol-7-yletter, rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

3H-[1]Benzopyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, acetate (ester), (3R,3aS)-rel- (9CI) (CA INDEX NAME) 452319-53-8 CAPLUS Z Z

Relative stereochemistry. Double bond geometry as shown.

Z Z

452319-55-0 CAPLUS
3H-[L]Benzopgyrano(4,3-c)isoxazol-7-ol, 3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, acetate (ester), (3R,3aS)-rel-(9CI) (GA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452319-57-2 CAPLUS
3H-IllBenzogyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, propanoate (ester),
[3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Z Z

452319-59-4 CAPLUS
Acetic acid, methoxy., (3R,3aS)-3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-ylester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

<12/04/2007>

10/513699

S S

452319-61-8 CAPLUS Cyclopropancezboxylic acid, (3R,3aS)-3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-yl ester, rel-(9C1) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452319-63-0 CAPLUS
Propanoic acid, 2,2-dimethyl-, (3R,3aS)-3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl] 3H-[l]benzopyrano[4,3-c]isoxazol-7-yl ester, rel- (9Cl) (CA INDEX NAME) £ £

Relative stereochemistry. Double bond geometry as shown.

2 Z

452319-65-2 CAPLUS 2-Propenoic acid, (3,3aS)-3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-ylester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

<12/04/2007>

452319-67-4 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 7-(cyclopentyloxy)-3a,4-dihydro-3-[[4-[(2s)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9Cl) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

S S

Relative stereochemistry. Double bond geometry as shown.

Z Z

452319-71-0 CAPLUS Ethanamine, 2-[[(3R,3aS)-3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-yl]oxy]-N.N-dimethyl-, rel- (9Cl) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

<12/04/2007>

Erich Leese

10/513699

Z Z

452319-73-2 CAPLUS
3H-[1]Benzogyrano[4,3-c]1soxazol-7-ol, 3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, hydrogen sulfite (ester), (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452319-75-4 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazole, 8-chloro-3a,4-dihydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

Z Z

452319-77-6 CAPLUS
3H-[1]Benzopyranol, 43-c]isoxazole, 8-chloro-3a,4-dihydro-7-methoxy-3-[[4[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel[9C1) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

S S

452319-78-7 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazole, 8-bromo-3a,4-dihydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

C. P.

452319-80-1 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 8-bromo-3a,4-dihydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(+)- (9CI) (CA INDEX NAME)

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

Z Z

452319-81-2 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 8-bromo-3a,4-dihydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(-)- (9CI) (CA

Erich Leese <12/04/2007>

10/513699

INDEX NAME)

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

452319-83-4 CAPLUS
3H-[11 Benzopyrano[4,3-c]isoxazole, 7,8-difluoro-3a,4-dihydro-3-[[4-[(2E)-2-apthyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452319-85-6 CAPLUS

[1] Renzopyrano [4,3-c] isoxazole, 8-fluoro-3a,4-dihydro-7-methoxy-3-[[4-] [(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

452319-87-8 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-8-phenyl-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452319-89-0 CAPLUS helllBenzopyrano[4,3-c]isoxazole, 3a,4-dihydro-8-methyl-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) ₹ S

Relative stereochemistry. Double bond geometry as shown.

452119-91-4 CAPLUS
3H-[1] Benzopyrano (4,3-c)isoxazole, 8-fluoro-3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-7-(methylthio)-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Erich Leese

<12/04/2007>

10/513699

452319-93-6 CAPLUS

##11]Bencapyrano[4,3-c]isoxazole-7,8-diol, 3a,4-dihydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

●2 HCl

452319-95-8 CAPLUS
3H-[1] Benzopyrano[4,3-c] isoxazol-8-ol, 3a,4-dihydro-7-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) E 5

Relative stereochemistry. Double bond geometry as shown.

Erich Leese

452319-97-0 CAPLUS Z Z

3H-[1]Benzopyrano[4,3-c]isoxazol-8-ol, 3a,4-dihydro-7-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, acetate (ester), (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452319-99-2 CAPLUS Ethanol, 2-[(18.3.436)-3a,4-dihydro-7-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyranol4,3-c]iscxazol-8-ylloxyl-, acetate (ester), rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452120-01-3 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) 2 3

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

452320-03-5 CAPLUS
4-11Bencopyranol(4,3-c)isoxazol-7-ol, 3-[[4-[(2E)-3-(3-fluorophenyl)-2-methyl-2-propenyl]-1-piperazinyl]methyl]-3a,4-dihydro-8-methoxy-, acetate (ester), (3R,38)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

452320-06-8 CAPLUS
#1.118enzopyzano! (4):2:13soxazol-7-01, 3-[[4-[(2E)-3-(4-fluorophenyl)-2-methyl-2-pyropenyl]-1-piperazinyl]methyl]-3a,4-dihydro-8-methoxy-, acetate (ester), (3R,3as)-rel-(9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452120-07-9 CAPLUS

1-11]Renzopyranol (4.3-c)1soxazol-7-ol, 3-[[4-[[2E]-3-(4-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-1a.4-dihydro-8-methoxy-, acetate (ester), [3R, JaS]-rel- (9CI) (CA INDEX NAME) **3** 5

Relative stereochemistry. Double bond geometry as shown.

S S

452320-09-1 CAPLUS
3H-[11Benzopyrano[4, 3-c]isoxazole, 7-(cyclopropylmethoxy)-3-[[4-[(2E)-3-(4-Eluorophenyl]-2-methyl-2-propenyl]-1-piperazinyl]methyl}-3a,4-dihydro-8-methoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452320-11-5 CAPLUS

Cyclopropanecarboxylic acid, (3R,3aS)-3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-1-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-yl ester, rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

<12/04/2007>

Erich Leese

10/513699

3-Pyridinecarboxylic acid, (3R,3aS)-3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-yl ester, rel- (9CI) (CA INDEX NAME) 452320-13-7 CAPLUS S S

Relative stereochemistry. Double bond geometry as shown.

S. S.

452320-15-9 CAPLUS carbair. (3R,3aS)-3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-carbamic acid, ethyl., (1-piperazlinyl]methyl]-3-phenyl-2-propenyl]-1-piperazlinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-yl ester, real. (9C1) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452320-17-1 CAPLUS Ethanamine, 2-[[48,3a8]-3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-yl]oxyl-N,N-dimethyl-, rel- [9CI] (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

<12/04/2007>

452320-19-3 CAPLUS
3H-[1]Benzogyrano[4,3-c]isoxazole, 3a,4-dihydro-8-methoxy-7-[(2-methoxyehoxy)methoxy]-3-[[4-[(2B)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9Cl) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452320-21-7 CAPLUS cacid, (3x, JaS)-3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-Phenyl-2-properyl]-1-piperainyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7-yl methyl ester, rel- (9Cl) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

3 3

452320-23-9 CAPLUS

##11]Benzopyrano(4,3-c]isoxazole, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-7-(2-propenyloxy)-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

<12/04/2007>

Erich Leese

Relative stereochemistry. Double bond geometry as shown.

Z Z

452320-25-1 CAPLUS
Ethanol, 2-[(13R, 348)-38, 4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-3H-[1]benzopyrano[4,3-c]isoxazol-7yl]oxyl-, dihydrochloride, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

●2 HCl

452320-27-3 CAPLUS
3H-[1]BenzOpyrano[4,3-c]iscxazol-7-ol, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, methanesulfonate (ester), (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452320-29-5 CAPLUS

3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452320-31-9 CAPLUS
41-[1]encopyrano[4,3-c]isoxazole, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-7-phenyl-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) **3** 8

Relative stereochemistry. Double bond geometry as shown.

Z Z

452320-34-2 CAPLUS
34-11Barapyranol4,3-clisoxazole, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)-2-ab-11)-2-phenyl-2-propenyl]-1-piperazinyl]methyl]-7-(methyl-thio)-, (3R,3aS)-rel-, mono(trifluoroacetate) (9Cl) (CA INDEX NAME)

<12/04/2007>

10/513699

ξ

CRN 452320-33-1 CMF C27 H33 N3 O3 S

Relative stereochemistry. Double bond geometry as shown.

£

76-05-1 C2 H F3 O2 CRN

C-C02H

452320-36-4 CADLUS Isoxazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452320-38-6 CAPLUS Isoxazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) ¥ 5

<12/04/2007>

Relative stereochemistry. Double bond geometry as shown.

Z Z

452320-40-0 CAPLUS Isoxazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452320-42-2 CAPLUS ISOSAZAOL(4,3-C)quinoline, 3-[[4-[(2E)-3-(3-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3,3a,4,5-tetrahydro-7,8-dimethoxy-, (3R,3aR)-rel-(9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Z 3

452320-44-4 CAPLUS Isoxazolo[4,3-c]quinoline, 3-[[4-[(2E)-3-(3-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3,3a,4,5-tetrahydro-7,8-dimethoxy-, (3R,3aR)-rel-(+)-

<12/04/2007>

10/513699

(9CI) (CA INDEX NAME)

Rotation (+). Absolute stereochemistry unknown. Double bond geometry as shown.

452320-46-6 CAPLUS 1203a20.0[4,3-c[quinoline, 3-[[4-[(2E)-3-(3-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3,3a,4,5-tetrahydro-7,8-dimethoxy-, (3R,3aR)-rel-(-)-(9Cl) (CA INDEX NAME) Z Z

Rotation (-). Absolute stereochemistry unknown. Double bond geometry as shown.

452320-48-8 CAPLUS
Isoxazolo[4,3-c[quinolline, 3-[[4-[(2E)-3-(4-fluorophenyl)-2-propenyl]-1-piperazinyl]methyl]-3,3a,4,5-tetrahydro-7,8-dimethoxy-, (3R,3aR)-rel-[9CI] (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Erich Leese

15oxazolo[4,3-0]quinoline, 3,3a,4,5-tetrahydro-7-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) 452320-50-2 CAPLUS Z Z

Relative stereochemistry. Double bond geometry as shown.

452320-52-4 CAPLUS Stokazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-5-methyl-3-[[4-[(2E)-3-phsnyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452320-54-6 CAPLUS CAPLUS CAPLUS 3,3a,4,5-tetrahydro-7,8-dimethoxy-5-methyl-3-[[4-[(28)-2-methyl-3-propenyl]-1-piperazinyl]methyl]-, (3R.3aR)-rel- (2G.] (CA INDEX NAME) S 53

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

452320-56-8 CAPLUS Isoxazolo[4].3-Gjuinoline, 3,3a,4,5-tetrahydro-7-methoxy-3-[[4-[(2E)-2-methyl-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-5-(phenylmethyl)-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) 2 S

Relative stereochemistry. Double bond geometry as shown.

452320-58-0 CAPLUS Isoxasolo(4,0-clquinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-5-(trifluoroacetyl)-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452320-60-4 CAPLUS Sozazalold,4,3-Clquinoline, 3,3a,4,5-tetrahydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-5-(trifluoroacetyl)-, (3K,3aR)-rel- (9CI) (CA INDEX NAME) 2 2

Relative stereochemistry. Double bond geometry as shown.

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Z Z

452320-62-6 CAPLUS Isoxazolo[4,3-c]quinoline, 5-acetyl-3,3a,4,5-tetrahydro-3-[[4-[[2E]-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452320-64-8 CAPLUS Isosazolo[4,3-c]quinoline-5(3H)-carboxylic acid, 3a,4-dihydro-3-[[4-[(2E)-3-phenyl-2-propenyl-1-piperazinyl]methyl]-, ethyl ester, (3R,3aR)-rel-9CI) (CA INDEX NAME) Z 23

Relative stereochemistry. Double bond geometry as shown.

452320-66-0 CAPLUS Soxazolo[4,3-c]quinoline-5(3H)-carboxamide, N-ethyl-3a,4-dlhydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

452320-68-2 CAPLUS
Naphrh[1,2-1]isoxazole, 3,3a,4,5-tetrahydro-7-methoxy-3-[[4-[(2E)-3-phenyl2-propenyl-1-piperazinyl]methyl]-, dihydrochloride, (3R,3aR)-rel- (9CI)
(CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

●2 HCl

452320-70-6 CAPLUS
Naphth[1,2-c]isoxazole, 3,3a,4,5-tetrahydro-3-[[4-[(2E)-3-phenyl-2-propenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

452330-72-8 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-3-[[4-(phenylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Relative stereochemistry.

Z Z

452320-74-0 CAPLUS Z.

<12/04/2007>

CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-3-[[4-(3-phenylpropyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9Cl) (CA INDEX NAME)

Relative stereochemistry.

Z Z

Relative stereochemistry.

●2 HCl

S S

452320-78-4 CAPLUS

[1] Benzopyrano [4, 3-c] isoxazole, 3a, 4-dihydro-7-[(2-methoxyethoxy) methoxyethoxy] -3-[[4.(2-naphthalenylmethyl)-1-piperazinyl]methyl] - (3R, 3aS)-rel- (9CL) (CA INDEX NAME)

Relative stereochemistry.

PAGE 1-A

Erich Leese <12/04/2007>

10/513699

PAGE 1-B

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(CA INDEX 3H-[1]Benzopyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) 452320-80-8 CAPLUS 2 Z

Relative stereochemistry.

452320-82-0 CAPLUS
4-11]Bencopyrano[4,3-c]isoxazole, 7-(cyclopropylmethoxy)-3a,4-dihydro-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry.

452320-84-2 CAPLUS
3H-[1] Benzopyrano[4,3-c] isoxazole, 7-(difluoromethoxy)-3a,4-dihydro-3-[[4[2-naphthalenylmethyl)-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME) Z Z

Relative stereochemistry.

Z Z

452320-86-4 CAPLUS

#11]Renzopyranol(4,3-c)isoxazole, 3a,4-dihydro-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-7-[(tetrahydro-3-furanyl)oxy]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry

452320-88-6 CAPLUS
4-11.llencopyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-3-[[4-(2-naphthalenylmethyl]-1-piperazinyl]methyl]-, acetate (ester), (3R,3aS)-rel-(9CI) (CA INDEX NAME) S 25

Relative stereochemistry.

Z Z

Relative stereochemistry.

452320-92-2 CAPLUS 3H-[1]Benzopyrano[4,3-c]isoxazol-7-ol, 3a,4-dihydro-8-methoxy-3-[[4-[(2E)-3-phenyl-2-butenyl]-1-piperazinyl]methyl]-, acetate (ester), (3R,3aS)-rel-2 Z

<12/04/2007>

Erich Leese

10/513699

(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452320-94-4 CAPLUS

4F21320-94-4 (A) -2 | Goxazole, 3-[[4-[(2E)-3-(2-fluorophenyl)-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3E,3aS)-rel-(9Cl) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452320-96-6 CAPLUS
#1.LlBencopyrano(4,3-c)isoxazole, 3-[[4-[(2E)-3-(3-fluorophenyl)-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

3 3

452320-98-8 CAPLUS
3H-[1]Bencopyrano[4,3-c]isoxazole, 3-[[4-[(2E)-3-(4-fluorophenyl)-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Z Z

452121-00-5 CAPLUS 1-[1]8encopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(4-methoxyphenyl)-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Z Z

452121-02-7 CAPLUS
4-11]sencepyranol(4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-[3-[1-(trifluoromethyl)phenyl]-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

452321-04-9 CAPLUS
3H-[1]Benzogpyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3]-[4-(trifluoromethyl)phenyl]-2-butenyl]-1-piperazinyl]methyl]-,
(3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452321-06-1 CAPLUS

[1] Rencopyrano (4,3-c) isoxazole, 3-[[4-[(2E)-3-(2-chlorophenyl)-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9Cl) (CA INDEX NAME) **3** 3

Relative stereochemistry. Double bond geometry as shown.

452321-08-3 CAPLUS

##1.[1]Rencopyrano(4,3-c]isoxazole, 3-[[4-[(2E)-3-(3-chlorophenyl)-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3E,3aS)-rel-(9Cl) (CA INDEX NAME) S S

Relative stereochemistry. Double bond geometry as shown.

452321-10-7 CAPLUS

4F.113encopyranol(4.3-c]isoxazole, 3-[[4-[(2E)-3-(4-chlorophenyl)-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3E,3aS)-rel-(9Cl) (CA INDEX NAME) **3** 3

Relative stereochemistry. Double bond geometry as shown.

452321-12-9 CAPLUS
4F1]8enzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(4-methylphenyl)-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

452321-14-1 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(2-methoxyphenyl)-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

452321-16-3 CAPLUS
3H-[1]Benzopyyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(3-methoxyphenyl)-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) 2 3

Relative stereochemistry. Double bond geometry as shown.

452321-19-6 CAPLUS
3H-[11Benzogyrano[4,3-c]isoxazole, 3-[[4-[(22)-3-f]uoro-3-phenyl-2-propertyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-,dihydrochloride, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

2 HCl

452321-21-0 CAPLUS
4F11]Rencapyrano(4,3-c)isoxazole, 3-[[4-[(2E)-3-(2,4-difluorophenyl)-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9CI) (CA INDEX NAME) 2 2

Relative stereochemistry. Double bond geometry as shown.

Z Z

452321-23-2 CAPLUS
452321-23-2 CAPLUS
4-11.Bencopyrano(4,3-c]isoxazole, 3-[[4-[(2E)-3-(3,4-difluorophenyl)-2-butenyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel-(9Cl) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

Erich Leese <12/04/2007>

10/513699

452321-25-4 CAPLUS
3H-[11Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3]-(2-pyridinyl)-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)
(CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

452321-27-6 CAPLUS

49-11:Renzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(3-pyridinyl)-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI)

(CA INDEX RAME) S S

Relative stereochemistry. Double bond geometry as shown.

452121-29-8 CAPLUS H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7,8-dimethoxy-3-[[4-[(2E)-3-(2-thieny])-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. Double bond geometry as shown.

<12/04/2007>

RN 45231-31-2 CAPLUS
CN 3H-[1]Berzopyzano[4,3-c]isoxazole, 3-[[4-[(2E)-3-(2-furanyl)-2-butenyl]-1piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA
INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

RN 452321-33-4 CAPLUS
CN 1soxsazolo(4,3-clquinoline, 3-[[4-[(4-chlorophenyl)methyl]-1piperazinyllmethyl]-3,3a,4,5-tetrahydro-7,8-dimethoxy-, (3R,3aR)-rel(9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452321-35-6 CAPLUS
CN ISOxazolo(4,-10-1)-1piperaziny]methyl]-1,3,3,4,5-tetrahydro-7,8-dimethoxy-, (3R,3aR)-rel(9CI) (CA INDEX NAME)

<12/04/2007> Erich

10/513699

Relative stereochemistry.

RN 452321-37-8 CAPLUS
CN Isoxazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[[4-[(2E)-3-phenyl-2-butenyl]-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry. Double bond geometry as shown.

RN 452321-39-0 CAPLUS
CN Isoxazolo(4,3-c)quinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452321-41-4 CAPLUS
CN Isoxazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[[4-(6quinolinylmethyl)-1-piperazinyl]methyl]-, (3R,3aR)-rel- (9CI) (CA INDEX

Erich Leese

NAME)

Relative stereochemistry.

RN 452321-43-6 CAPLUS
CN Isoxacolo(4,-2-ciquinoline, 3-[[4-[(5-fluoro-2-naphthalenyl)methyl]-1piperazinyl]methyl]-3,3a,4,5-tetrahydro-7,8-dimethoxy-, (3R,3aR)-rel(CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452321-45-8 CAPLUS
CN Isoxazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-7,8-dimethoxy-3-[[4-[2-(2-napkhalenyloxy)ethy1]-1-piperaziny1]methy1]-, (3R,3aR)-rel- (9CI) (CAINDEX NAME)

Relative stereochemistry.

RN 452321-47-0 CAPLUS

<12/04/2007> Erich Leese

10/513699

CN Isoxazolo[4,3-c]quinoline, 3,3a,4,5-tetrahydro-5-methyl-3-[[4-(2-naphthalenylmethyl)-1-piperazinyl]methyl]-, (3R,3aR)-rei- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452321-55-0 CAPLUS
CN 3H-[1]Benzopyrano[4,3-c]isoxazole, 3a,4-dihydro-7-methoxy-3-[2-[4-(2-naphthalenylmethyl)-1-piperazinyl]ethyl]-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

RN 452321-57-2 CAPLUS
CN 3H-[1] Benzopyrano[4,3-c] isoxazole-3-methanamine, 3a,4-dihydro-7,8-dimethoxy-N-methyl-N-[1-(2-naphthalenylmethyl)-4-piperidinyll-, (3R,3aS)-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

<12/04/2007> Erich Leese

452321-59-4 CAPLUS
Commanide, N-[[[3]x 3aS]-3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-1-y1]methyl]-N-[1-(2-naphthalenylmethyl)-4-piperidinyl]-, rel-(9CI) (CA INDEX NAME) Z Z

Relative stereochemistry.

1-Piperazinemethanamine, 4-[[(3R,3aS)-3a,4-dihydro-7,8-dimethoxy-3H-[1]benzopyrano[4,3-c]isoxazol-3-yl]methyl]-N-2-naphthalenyl-, rel- (9CI) (CA INDEX NAME) 452321-61-8 CAPLUS Z Z

Relative stereochemistry.

452934-93-9 CAPLUS

[1] Benzopyrano[4,3-c] isoxazole, 3-[[4-[(7-fluoro-2,3,4a,8a-tetrahydro-1,4-dioxino[2,3-b] pyridin-3-yl] methyl] -1-piperazinyl] methyl] -3a,4-dihydro-7,8-dimethoxy-, (3R,3as)-rel- (9CI) (CA INDEX NAME) Z Z

Relative stereochemistry. [§]

Erich Leese <12/04/2007>

10/513699

452934-94-0 CAPLUS
3H-[11.Benzograno[4, 3-C]isoxazole, 3-[[4-[[6-fluoro-3,4,4a,8a-tetrahydro-4-nethyl-3H-1,4-benzoxazin-2-yl)methyl]-1-piperazinyl]methyl]-3a,4-dihydro-7,8-dimethoxy-, (3R,3aS)-rel- (9CI) (CA INDEX NAME) S S

Relative stereochemistry.

THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT N REFERENCE COUNT:

=> d ibib abs hitstr 2-6

Preparation of novel 4,5-dihydroisoxazole derivatives and their use as pharmaceuticals for T cell-mediated L5 ANSWER 2 OF 6 CAPLUS COPYRIGHT 2007 ACS on STN ACCESSION NUMBER: 120:29:357 DOCUMENT NUMBER: Preparation of novel 4,5-diby diseases INVENTOR (S) :

Freyne, Eddy Jean Edgard; Andres-Gil, Jose Ignacio, Deroose, Frederik Dirk; Petit, Davy Petrus Franciscus Maria; Matesanz-Ballesteros, Maria Encarnacion; Alvarez Escobar, Rosa Maria danssen Pharmaceutica N.V., Belg. PCT Int. Appl., 108 pp.

PATENT ASSIGNEE(S): SOURCE:

LANGUAGE: FAMILY ACC. NUM. COUNT: PATENT INFORMATION: DOCUMENT TYPE:

APPLICATION NO. KIND DATE PATENT NO.

DATE

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The invention concerns title compds. I and their N-oxides, pharmaceutically acceptable addition salts, quaternary ammonium salts, and stereochem. isomeric forms [wherein m, n, p = 0 or 1; R1 = (un)substituted pyridinyl, pyridazinyl, pyriazinyl or phenyl; B = amide, ketone, or oxadiazole; D = (un)substituted aryl or heterocyclyl; Q = bond, CO, (un)substituted NH, CONH, CH2, CH(:CH2), C(:NH), SO, SO, 3-oxobutenyl, pyrazole, isoxazole, or thiazole nucleus; L = (un)substituted aryl or heterozayl; R2, R3 = H, halo, C1-6 alkyloxy, or (un)substituted C1-6 alkyloxy, and substituted C1-6 alkyloxy and alkyloxy activity and their medical use. The compds. show growth inhibitory activity ¥B

Erich Leese <12/04/2007>

10/513699

against T cell blasts and keratinocytes in vitro. The compds. are claimed for use in the treatment of prevention of rheumatic, arthritic, and inflammatory diseases, psoriasis, T cell leukemia, transplant rejection, and graft-vs.-host disease. For instance, base-catalyzed cycloaddn. of N-hydroxy-3-pyridinecarboximidoyl chloride with Me 2-propenoate gave 98* Me 4,5-dihydro-3-(3-pyridinyl)-5-isoxazolecarboxylate, which was amidated with (4-aminophenyl)phenylmethanone to give 58* title compound II. At a concentration of 10-6 M, II gave 81* inhibition of T cell blast formation in human whole blood.

264606-57-7

H

RL: RCT (Reactant); RACT (Reactant or reagent)
(starting material; preparation of dihydroisoxazole derivs. as
antiproliferatives and immunomodulators)
264606-57-7 CAPLUS
ISOXAZOLO[4, 3-c] quinoline-3-carboxylic acid, 3,3a,4,5-tetrahydro- (9CI)
(CA INDEX NAME) £ §

264606-16-8P HO₂C II

Strict BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (target compound; preparation of dihydroisoxazole derive. as antiproliferatives and immunomodulators) 264666.1es CAPUS.

Piperidine, 4-(4-fluorobenzoyl)-1-[[(3R,3aR)-3,3a,4,5-tetrahydroisoxazole(4,3-c)] (CA INDEX

Z Z

Relative stereochemistry.

THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT 9 REFERENCE COUNT:

JS COPYRIGHT 2007 ACS on STN 1997:431584 CAPLUS 127:149098 LE ANSWER 3 OF 6 CAPLUS
ACCESSION NUMBER: 19:
DOCUMENT NUMBER: 12:

A convenient synthesis of 3- and 3,4-substituted 4,5-dihydroisoxazole-5-acetic acids Eichinger, Rarl, Wokurek, Michael; Zauner, Bernd, Roscami, Mohammad Reza Institute of Organic Chemistry, Vienna University Technology, Vienna, A-1060, Austria 2733-2742 AUTHOR (S): TITLE:

oŧ CORPORATE SOURCE:

SOURCE:

CODEN: SYNCAV, ISSN: 0039-7911

PUBLISHER: DOCUMENT TYPE:

English CASREACT 127:149098 LANGUAGE: OTHER SOURCE(S): GI

The 4,5-dihydroisoxazole-5-acetic acids I [RI = Ph, Me, 4-CIC6H4, 4-PhC6H4, R2 = H, SPh, OPh, 4-CIC6H4, R12 = (CH2)10, 1,2.3,4-terrahydronaphth-1,2-diyl] were prepared from the ketoximes RIC(CH2R2):NOH, 2,2-dimethyl-5-methoxymethylene-1,3-dioxan-4,6-dione and 191247-115thum in yields from 35 to 79 %.

RI: SPN (Synthetic preparation); PREP (Preparation) (preparation of isoxazoleacetic acids)
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193267-4 2 H

Z Z

(CA INDEX NAME)

Relative stereochemistry.

THERE ARE 16 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT 16 REFERENCE COUNT:

CAPLUS COPYRIGHT 2007 ACS on STN
1955:228191
CAPLUS
122:81272
Nitrile oxide [3 + 2] cycloaddition: application to the synthesis of 6-substituted 3[2H]-pyridazinones and 6-substituted 4,5-dihydro-4-hydroxy-3[2H]-pyridazinones L5 ANSWER 4 OF 6 ACCESSION NUMBER: DOCUMENT NUMBER:

<12/04/2007>

Erich Leese

10/513699

The source of the preparation of 6-substituted 3(2H)-pyridazinones and efficient method for the preparation of 6-substituted 3(2H)-pyridazinones and 6-substituted 4,5-dihydro-4-hydroxy-3(2H)-pyridazinones starting from 3,5-disubstituted 4,5-dihydroisoxazoles is described. N-O bond cleavage of the isoxazoline ring promoted by molyhdenum hexacarbonyl or by catalyvic hydrogenation afforded the a-hydroxy f-keto esters RCOCH2CH(OH)COZET (I. R = Me, Bu, 2-, 4-pyridyl, 4-HOC6H4) which were converted into 6-substituted 4,5-dihydrox-4-hydroxy-3(HH)-pyridazinones for 6-substituted 3(2H)-pyridazinones on treatment with hydrazine hydrate at room temperature or reflux in high yield starting from I. An intramol, version of this methodol has been developed to prepare the known antiulcer Baraldi, P. G.; Bigoni, A.; Cacciari, B.; Caldari, C.; Manfredini, S.; Spalluto, G. Dipartimento di Scienze Parmaceutiche, Univ. di Perrara, Ferrara, I-44100, Iraly Synthesis (1994), (11), 1158-62 (CDBN: SYNTBF, ISSN: 0039-7881 tricyclic 5H-[1]-benzopyrano[4,3-c]pyridazin-3(2H)-one. Journal DOCUMENT TYPE: LANGUAGE: OTHER SOURCE(S): CORPORATE SOURCE: AUTHOR (S): PUBLISHER: SOURCE:

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT Ħ

(Reactant or reagent)
(intrile oxide [3 + 2] cycloaddn. to pyridazinones)
160427-31-6 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole-3-carboxylic acid, 3a,4-dihydro-, methylester (9CI) (CA INDEX NAME) Z Z

LE ANSWER 5 OF 6 CAPLUS COPYRIGHT 2007 ACS on STN ACCESSION NUMBER: 1995:94454 CAPLUS DOCUMENT NUMBER: 123:111970

TITLE:

Pyridazin-3(2H)-ones via A2-isoxazoline intermediates: synthetic studies Baraldi, Pier Giovanni; Spalluto, Giampiero; Manfredini, Stefano; Simoni, Daniele Dipartimento di Scienze Parmaceutiche, Universita di Perrara, Ferrara, Italy Acta Chimica Slovenica (1994), 41(2), 149-72 Journal AUTHOR (S):

CORPORATE SOURCE:

SOURCE:

DOCUMENT TYPE: LANGUAGE: GI

An efficient method for the preparation of 6-substituted-4,5-dihydro-3(2H)pyridazinones I (R = alkyl, benzyl, etc.) was described. The synthetic strategy is based on d2-isoxazolines chemical which were unmasked by N-o bond cleavage and cyclized to the target compound Utilizing the same approach was possible to obtain both 6-substituted-3(2H)-pyridazinones and 6-substituted-4-hydroxy-4,5-dihydro-3(2H)-pyridazinones and 6-substituted-4-hydroxy-4,5-dihydro-3(2H)-pyridazinones and 6-substituted-4-hydroxy-4,5-dihydro-3(2H)-pyridazinones and 6-substituted-4-hydroxy-4,5-dihydro-3(2H)-pyridazinone also extended to a C-nucleoside starting from P-ribofuranosylnitromethane. Moreover, an intramol. version of this method. has been developed to prepare a known antiulor tricyclic corresponding 3-(1-naphthyl)propionic acid Et ester derivs. was also æ

reported. 160427-31-6P RL: RCT (Reactant), SPN (Synthetic preparation); PREP (Preparation), RACT (Reactant or reagent)

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(preparation of benzpyranopyridazinone)

160427-31-6 CAPLUS
3H-[1]Benzopyrano[4,3-c]isoxazole-3-carboxylic acid, 3a,4-dihydro-, methylester (9CI) (CA INDEX NAME) Z Z

A general synthetic route to fused furans. Total synthesis of (+) pallescensin A Shishido, Kozo; Umimoto, Koji; Ouchi, Mikiko; Irie, Oamu; Omodani, Tomoki; Takata, Takeshi; Shibuya, Masayuki L5 ANSWER 6 OF 6 CAPLUS COPYRIGHT 2007 ACS on STN ACCESSION NUMBER: 1993:234262 CAPLUS , DOCUMENT NUMBER: 118:234262 AUTHOR (S):

Inst. Med. Resour., Univ. Tokushima, Tokushima, 770, Japan CORPORATE SOURCE:

Journal of Chemical Research, Synopses (1993), (2), 58-9 CODEN: JRPSDC, ISSN: 0308-2342

SOURCE:

DOCUMENT TYPE: LANGUAGE: GI

Erich Leese

<12/04/2007>

. 10/513699

AB A general and facile synthetic route to fused furans has been developed. The key step of the transformation involves the intramol. [13-2] dipolar cycloaddm. reaction of nitrile oxides which were generated in situ from the corresponding oxima acetaes. Reductive hydrolysis of the resulting dihydroisoazacles followed by alkaline hydrolysis provided \$\beta\$. \$\frac{1}{2}\$ dihydroxy ketones which were immediately treated with a catalytic amount of p-tolenesulfonic acid to afford the fused furans I (R = Me, R = H; R = Me). Alternatively, the alcs., derived by hydrolysis of the dihydroisoazacles, were submitted to a sequential reductive hydrolysis of treatment to provide I. Addnl. dihydroisoazacle alcs., prepared from com. available phthalide, were similarly treated to give the tricyclic fused furans II (R2 = H, Me, R3 = H) in reasonable yields. The methodol. developed here has been successfully applied to a total synthesis of (+)-pallescensin A (III) starting with (+)-Wieland-Wiescher ketone.

IT 147378-13-0P 14751-14-6P PB H

(Reactant or reagent)

(preparation and deacetylation of)
17378-13-0 (ARDUS
Naphth[1,2-c]isoxazole-3-methanol, 3,3a,4,5-tetrahydro-a-methyl-,
acetate (estet), [34(8*),3ad]- (9CI) (CA INDEX NAME) £ 5

147511-14-6 CAPLUS Naphth[1,2-c]isoxazole-3-methanol, 3,3a,4,5-tetrahydro- α -methyl-, acetate (ester), [3 $\alpha(R^*)$,3a α]- (9CI) (CA INDEX NAME) Z Z

147378-22-1P 147511-12-4P

<12/04/2007>

Relative stereochemistry.

RN 147511-12-4 CAPLUS CN Nabhthil.2-clisoxazole-3-methanol, 3,3a,4,5-tetrahydro-α-methyl-, [3α(R*),3aα]- (9CI) (CA INDEX NAME)

Relative stereochemistry.

Erich Leese